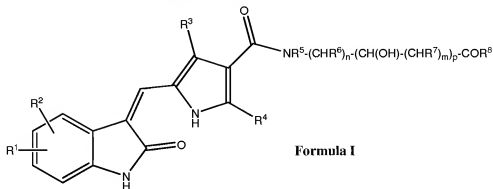


IN THE CLAIMS

1. (Original) A compound represented by Formula (I):



wherein:

R^1 is selected from the group consisting of hydrogen, halo, (C1-C6) alkyl, (C3-C8) cycloalkyl, (C1-C6) haloalkyl, hydroxy, (C1-C6) alkoxy, amino, (C1-C6) alkylamino, amide, sulfonamide, cyano, substituted or unsubstituted (C6-C10) aryl;

R^2 is selected from the group consisting of hydrogen, halo, (C1-C6) alkyl, (C3-C8) cycloalkyl, (C1-C6) haloalkyl, hydroxy, (C1-C6) alkoxy, (C2-C8) alkoxyalkyl, amino, (C1-C6) alkylamino, (C6-C10) arylamino;

R^3 is selected from the group consisting of hydrogen, (C1-C6) alkyl, (C6-C10) aryl, (C5-C10) heteroaryl, and amide;

R^4 , R^5 and R^6 are independently selected from the group consisting of hydrogen and (C1-C6) alkyl;

each R^7 is independently selected from the group consisting of hydrogen, (C1-C6) alkyl and hydroxyl;

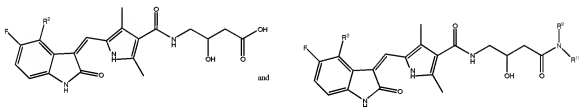
R^8 is selected from the group consisting of hydroxy, (C1-C6) O-alkyl, (C3-C8) O-cycloalkyl, and NR^9R^{10} ; where R^9 and R^{10} are independently selected from the group consisting of hydrogen, (C1-C6) alkyl, (C1-C6) hydroalkyl, (C1-C6) dihydroxyalkyl, (C1-C6) alkoxy, (C1-C6) alkyl carboxylic acid, (C1-C6) alkyl phosphoric acid, (C1-C6) alkyl sulfuric acid, (C1-C6) hydroxyalkyl carboxylic acid, (C1-C6) alkyl amide, (C3-C8) cycloalkyl, (C5-C8) heterocycloalkyl, (C6-C8) aryl, (C5-C8) heteroaryl, (C3-C8) cycloalkyl carboxylic acid, or R^9

and R^{10} together with N forms a (C5-C8) heterocyclic ring either unsubstituted or substituted with one or more hydroxyls, ketones, ethers, and carboxylic acids; and

n and m are independently 0, 1, 2, or 3; p is 1, 2, or 3;

or, a pharmaceutically acceptable salt, its tautomer, a pharmaceutically acceptable salt of its tautomer, or a prodrug thereof.

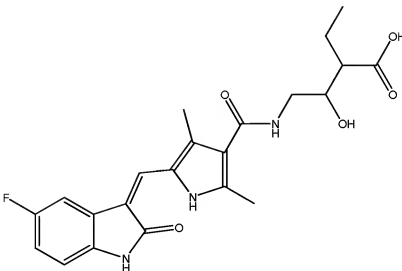
2. (Original) The compound, salt, tautomer, or prodrug according to claim 1 selected from the group represented by the following structures:



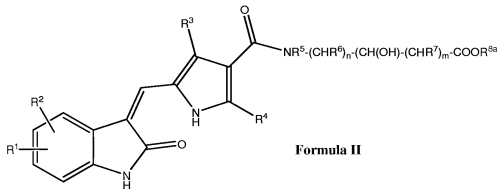
wherein

n R^2 is selected from the group consisting of hydrogen and fluoro.

3. (Original) The compound, salt, tautomer, or prodrug according to claim 1 represented by the following structure:



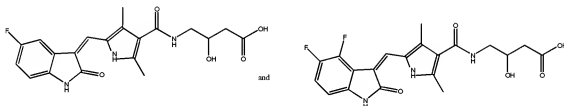
4. (Original) The compound, salt, tautomer, or prodrug according to claim 1 represented by Formula (II):



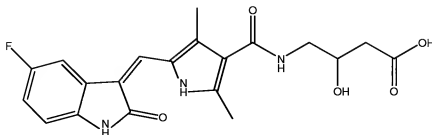
wherein R^{8a} is selected from the group consisting of hydrogen, (C1-C6) alkyl, and (C3-C8) cycloalkyl.

5. (Original) The compound, salt, tautomer, or prodrug according to claim 4, wherein:
 R^1 and R^2 are independently selected from the group consisting of hydrogen and fluoro;
 R^3 and R^4 are methyl;
 R^5 , R^6 , R^7 and R^{8a} are hydrogen; and
 n and m are independently 0, 1, or 2.

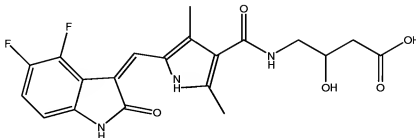
6. (Original) The compound, salt, tautomer, or prodrug according to claim 5 selected from the group consisting of:



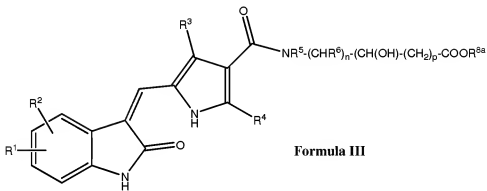
7. (Original) The compound, salt, tautomer, or prodrug according to claim 5 represented by the following structure:



8. (Original) The compound, salt, tautomer, or prodrug according to claim 5 represented by the following structure:



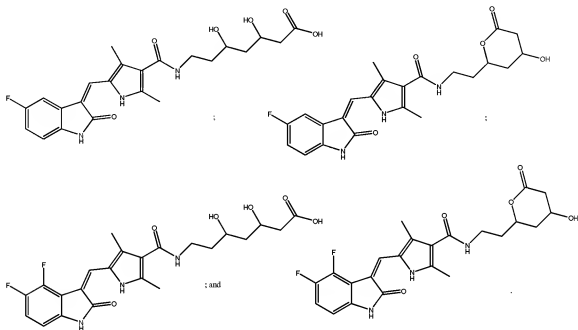
9. (Original) A compound, salt, tautomer, or prodrug according to claim 1 represented by Formula (III):



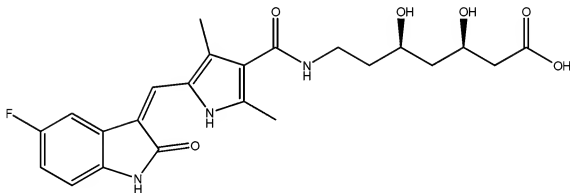
Formula III

wherein R^{8a} is selected from the group consisting of hydrogen, (C1-C6) alkyl, and (C3-C8) cycloalkyl.

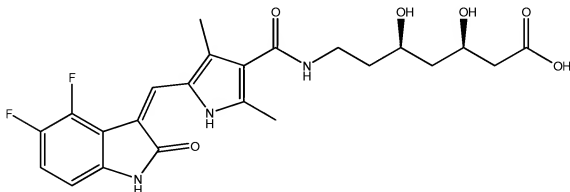
10. (Original) The compound, salt, tautomer, or prodrug according to claim 9, wherein:
 R^1 and R^2 are independently selected from the group consisting of hydrogen and fluoro;
 R^3 and R^4 are methyl;
 R^5 , R^6 , and R^{8a} are hydrogen; and
 n and p are independently 1, or 2.
11. (Original) The compound, salt, tautomer, or prodrug according to claim 10 selected from the group consisting of:



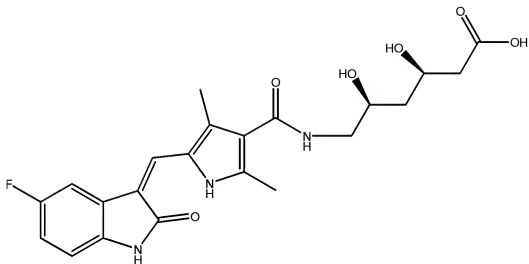
12. (Original) The compound, salt, tautomer, or prodrug according to claim 10 represented by the following structure:



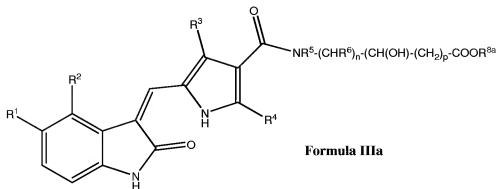
13. (Original) The compound, salt, tautomer, or prodrug according to claim 10 represented by the following structure:



14. (Original) The compound, salt, tautomer, or prodrug according to claim 10 represented by the following structure:



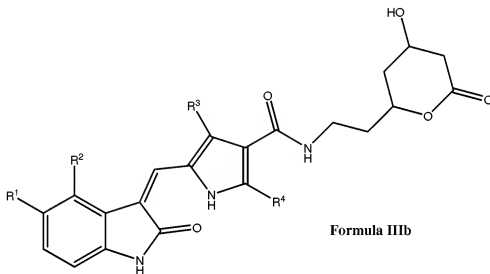
15. (Original) A compound, salt, tautomer, or prodrug according to claim 9 represented by Formula (IIIa):



wherein:

R^1 and R^2 are independently selected from the group consisting of hydrogen and fluoro;
 R^3 and R^4 are methyl;
 R^5 , R^6 , and R^{8a} are hydrogen; and
 n and p are 2.

16. (Original) A compound, salt, tautomer, or prodrug according to claim 15 represented by Formula (IIIb):

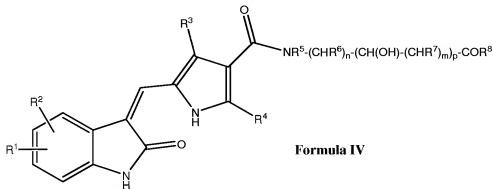


Formula IIIb

wherein:

R^1 and R^2 are independently selected from the group consisting of hydrogen and fluoro;
and
 R^3 and R^4 are methyl.

17. (Original) A compound, salt, tautomer, or prodrug according to claim 1 represented by Formula (IV):



wherein R^8 is NR^9R^{10} .

18. The compound, salt, tautomer, or prodrug of claim 17, wherein:

R^1 and R^2 are independently selected from the group consisting of hydrogen, halo, cyano;

R^3 , R^4 , R^5 and R^6 are independently hydrogen or (C1-C6) alkyl;

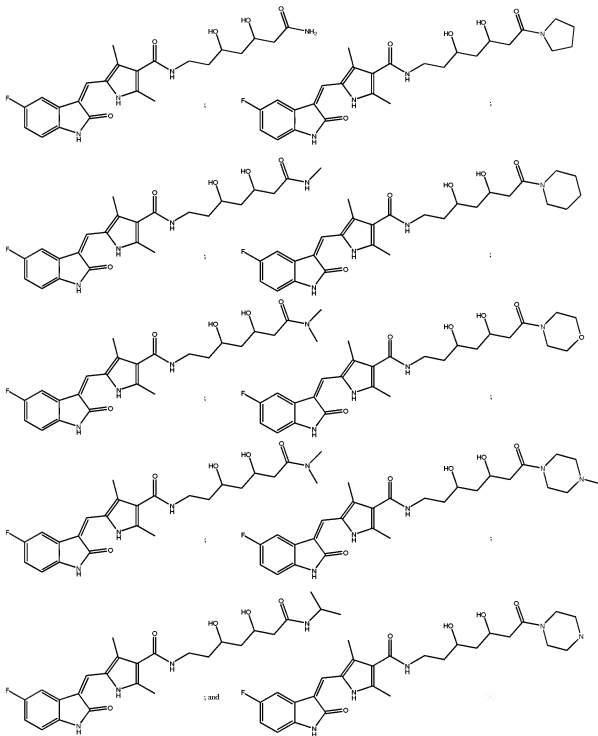
R^7 is hydrogen, or hydroxyl;

n , and p are independently 1, or 2;

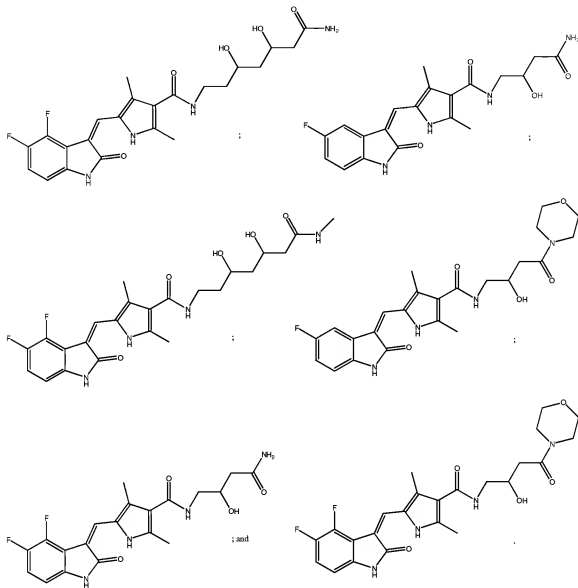
m is 0 or 1; and

R^9 and R^{10} are selected from the group consisting of hydrogen, (C1-C6) alkyl, (C1-C6) hydroxyalkyl, (C1-C6) dihydroxyalkyl, (C1-C6) alkoxy, (C1-C6) alkyl carboxylic acid, (C1-C6) alkyl phosphoric acid, (C1-C6) alkyl sulfuric acid, (C1-C6) hydroxyalkyl carboxylic acid, (C1-C6) alkyl amide, (C3-C8) cycloalkyl, (C5-C8) heterocycloalkyl, (C6-C8) aryl, (C5-C8) heteroaryl, (C3-C8) cycloalkyl carboxylic acid, or R^9 and R^{10} together with N forms a (C5-C8) heterocyclic ring either unsubstituted or substituted with one or more hydroxyls, ketones, ethers, and carboxylic acids.

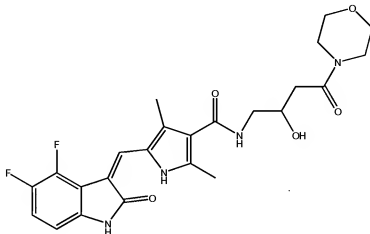
19. (Original) The compound, salt, tautomer, or prodrug according to claim 18 selected from the group represented by the following structures:



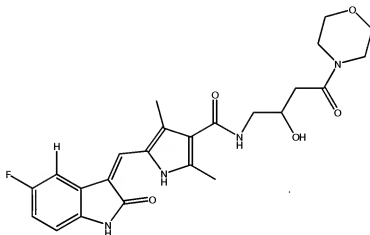
20. (Original) The compound, salt, tautomer, or prodrug according to claim 18 selected from the group represented by the following structures:



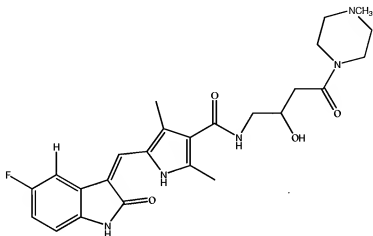
21. (Original) The compound, salt, tautomer, or prodrug according to claim 18 represented by the following structure:



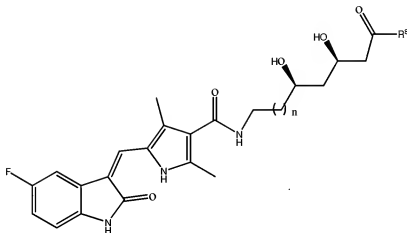
22. (Original) The compound, salt, tautomer, or prodrug according to claim 18 represented by the following structure:



23. (Original) The compound, salt, tautomer, or prodrug according to claim 18 represented by the following structure:

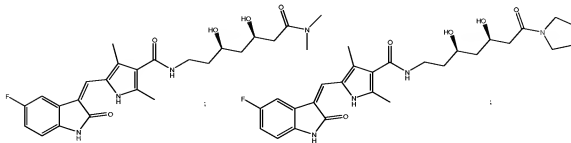


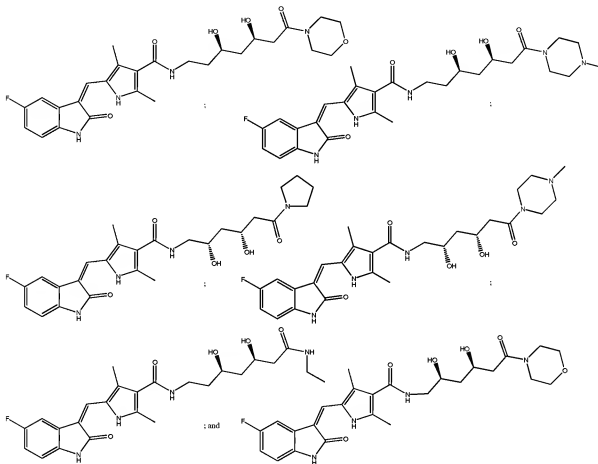
24. (Original) The compound, salt, tautomer, or prodrug according to claim 18 represented by the following structure:



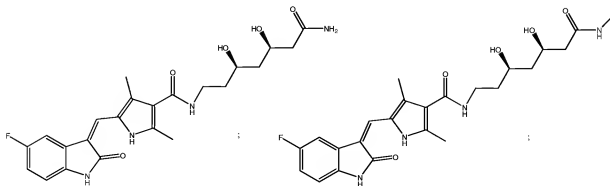
wherein n is 0, 1, or 2.

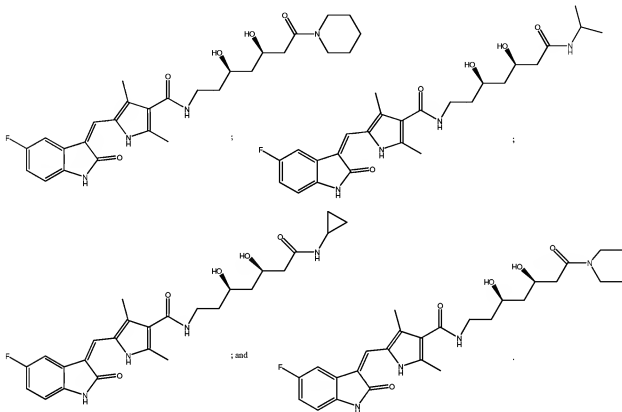
25. (Original) The compound, salt, tautomer, or prodrug according to claim 24 selected from the group represented by the following structures:



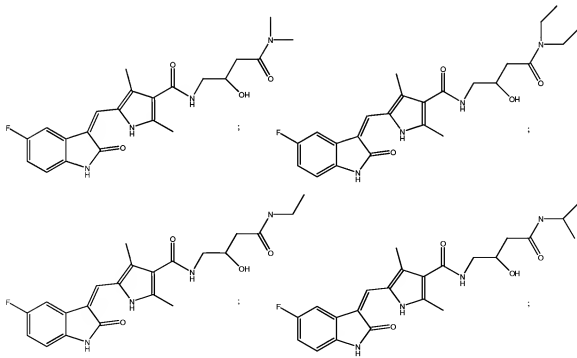


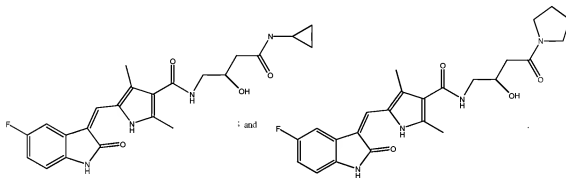
26. (Original) The compound, salt, tautomer, or prodrug according to claim 24 selected from the group represented by the following structures:



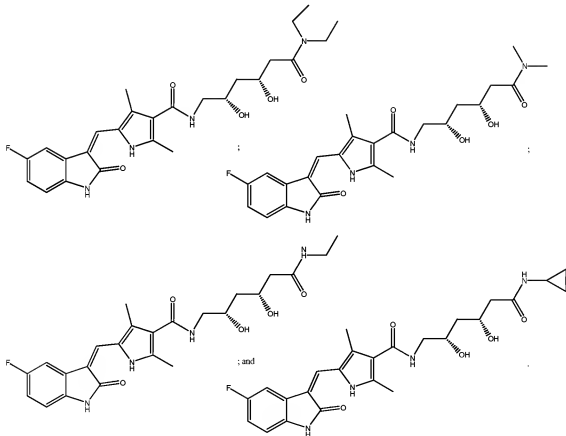


27. (Original) The compound, salt, tautomer, or prodrug according to claim 18 selected from the group represented by the following structures:

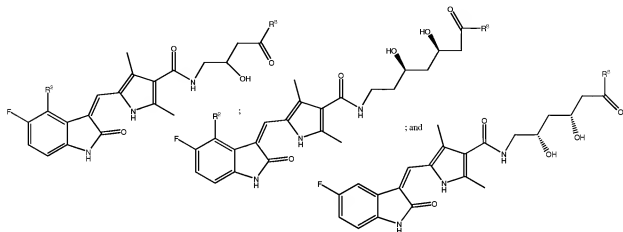




28. (Original) The compound, salt, tautomer, or prodrug according to claim 18 selected from the group represented by the following structures:



29. (Original) The compound, salt, tautomer, or prodrug according to claim 18 selected from the group represented by the following structures:

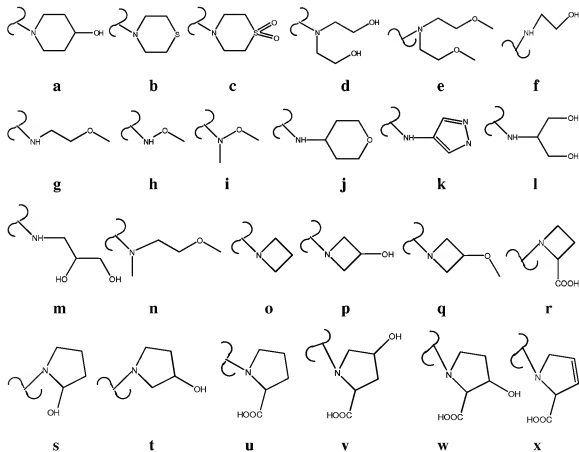


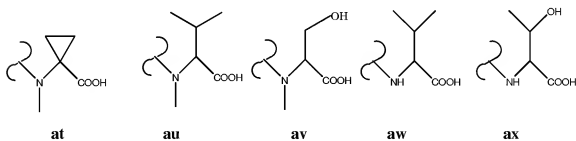
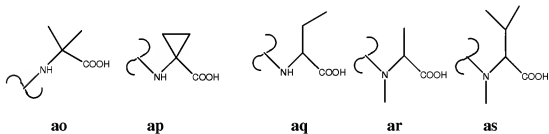
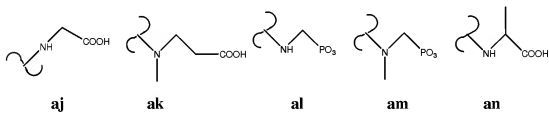
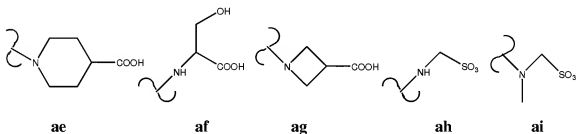
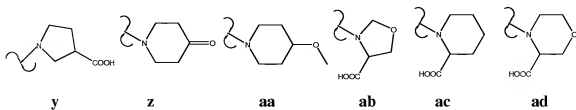
wherein:

R^2 is selected from the group consisting of hydrogen and fluoro; and

R^8 is selected from the group consisting of radicals represented by the following

structures:





30. (Canceled)

31.-32. (Canceled)